

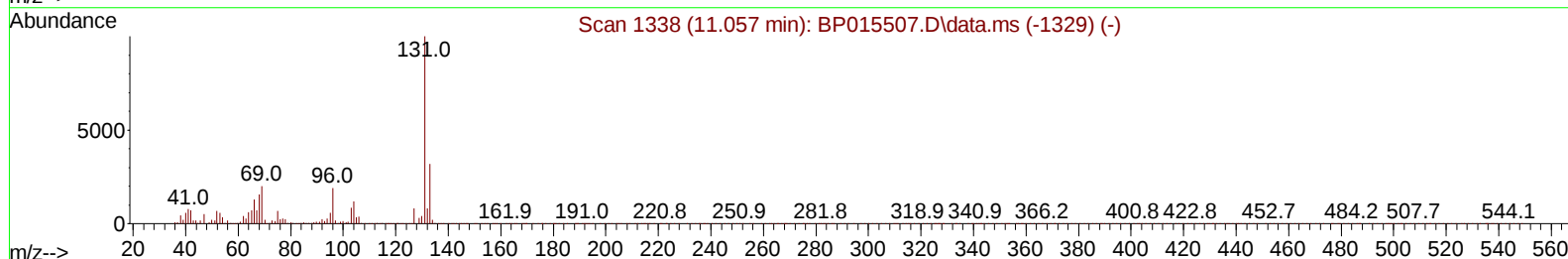
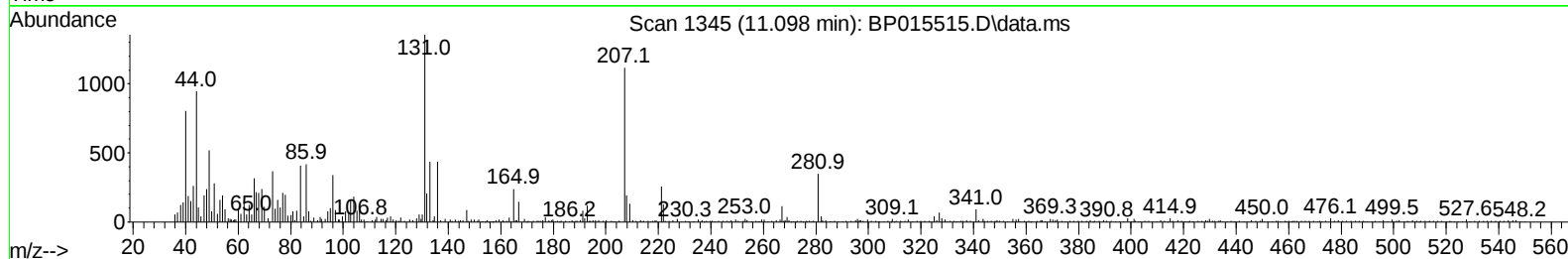
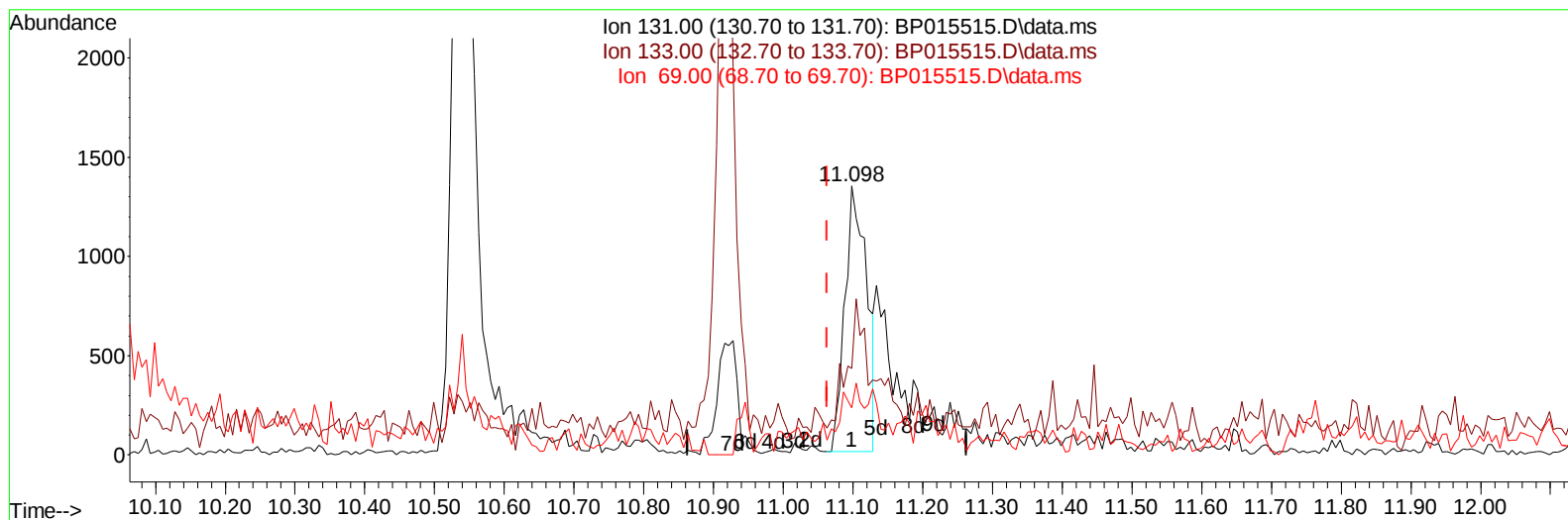
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015515.D
Acq On : 14 Jun 2023 13:48
Operator : MA/JU
Sample : 03098-10
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EZEP7

Manual IntegrationsAPPROVED

Quant Time: Jun 14 23:09:14 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 09 23:33:38 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 06/15/2023
Supervised By :mohammad ahmed 06/16/2023



TIC: BP015515.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.098min (+ 0.035) 0.19 ng/u1

response 2886

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	32.25
69.00	18.30	17.56
0.00	0.00	0.00

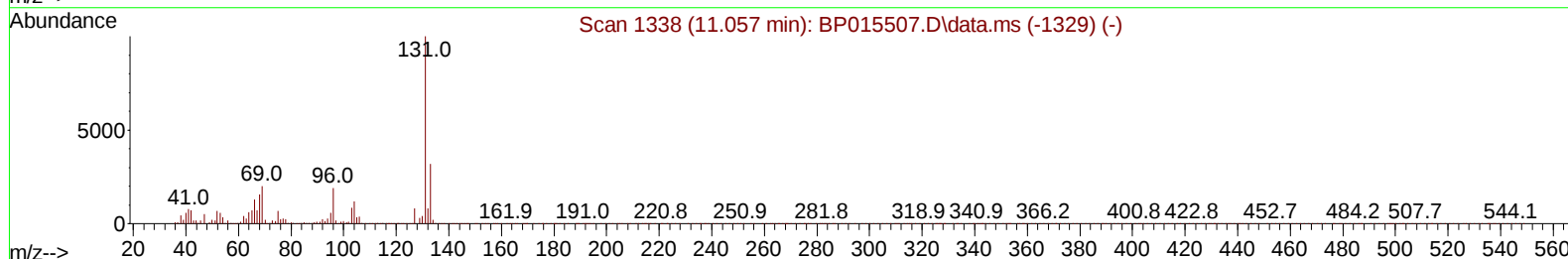
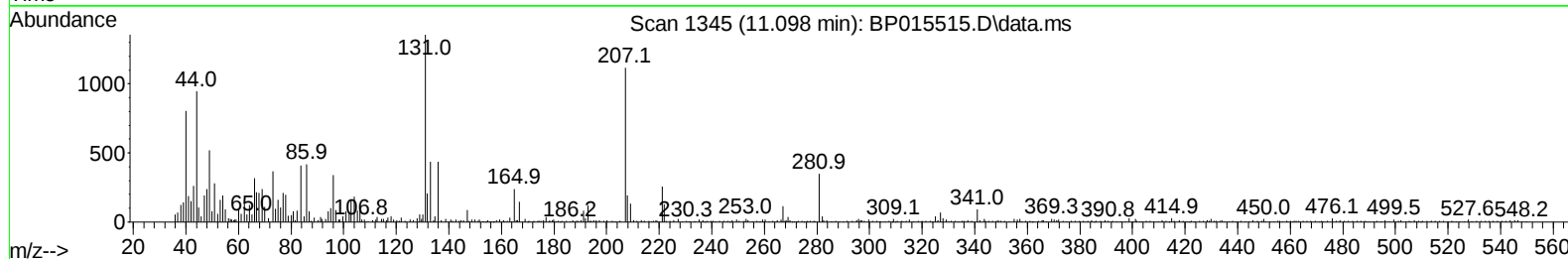
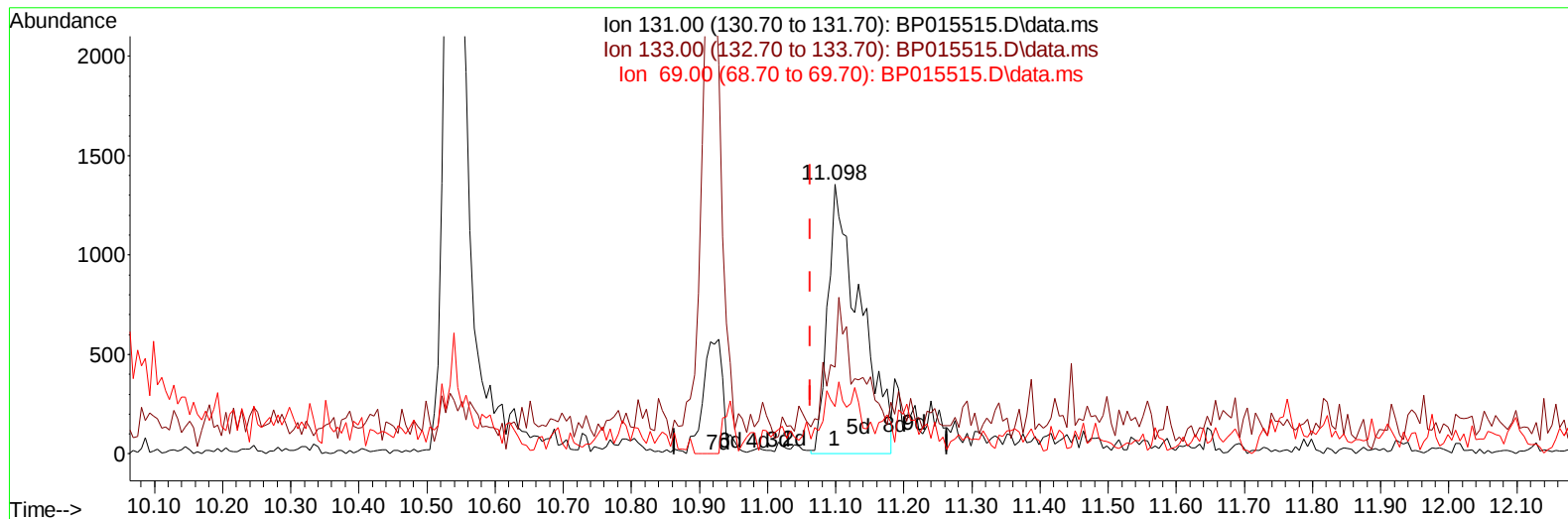
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TIC: BP015515.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.098min (+ 0.035) 0.29 ng/ul m

response 4458

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	32.25
69.00	18.30	17.56
0.00	0.00	0.00

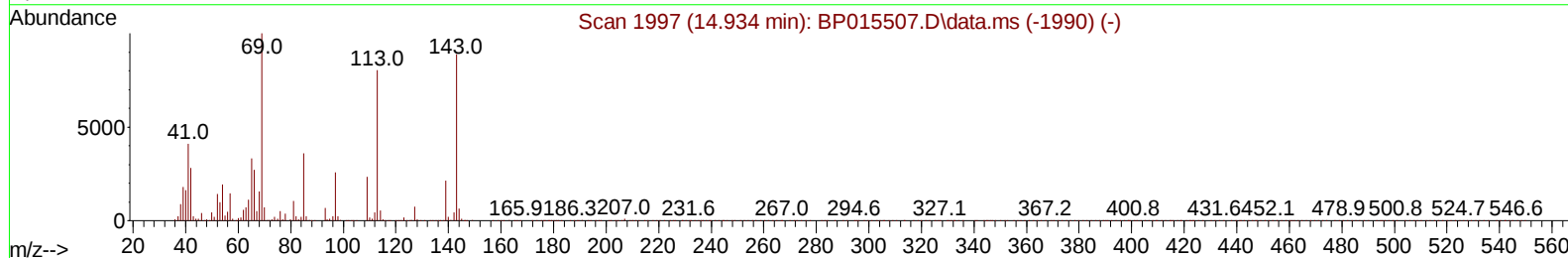
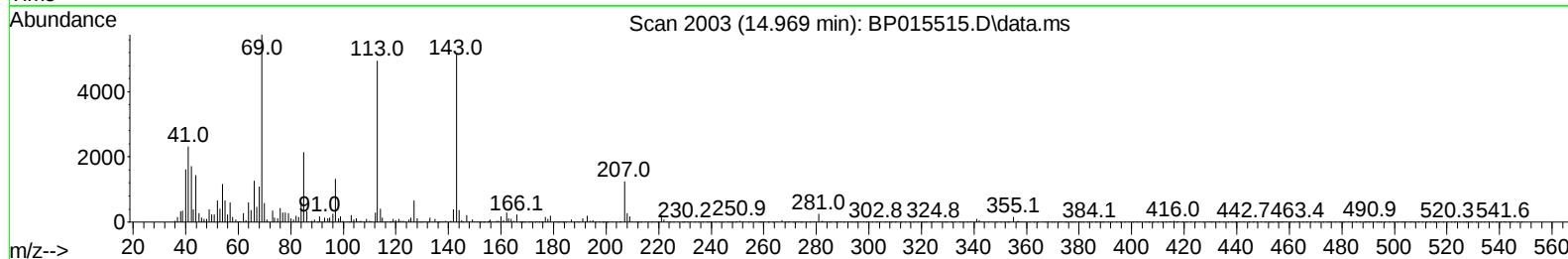
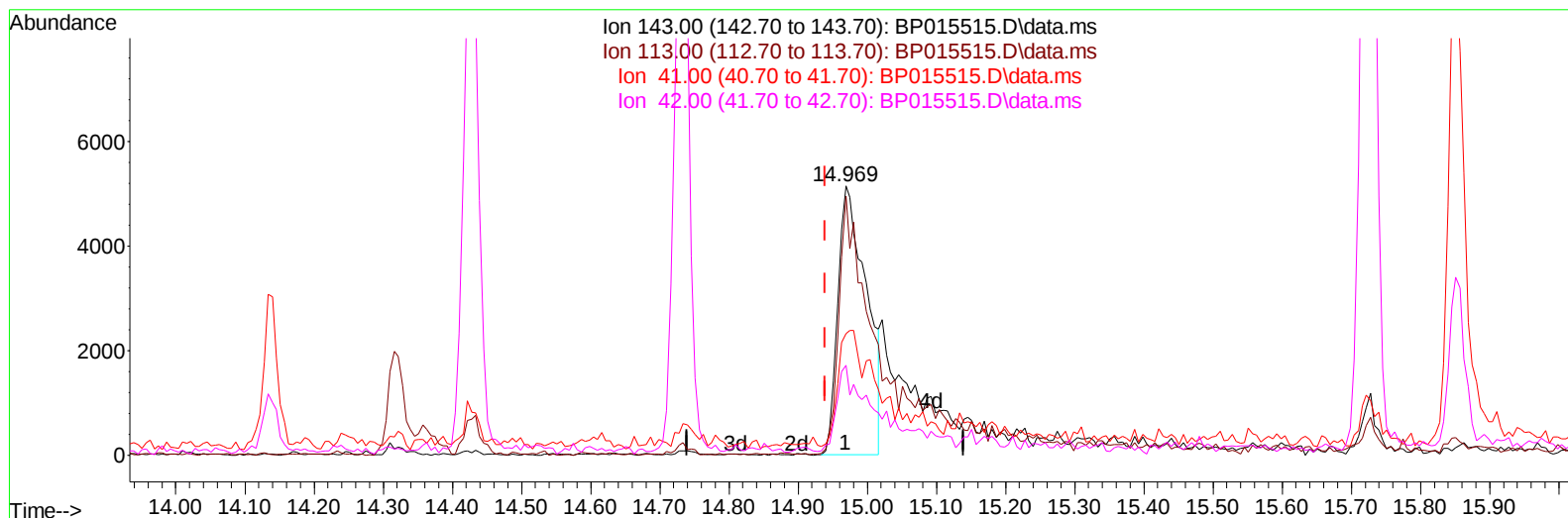
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Instrument :
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TIC: BP015515.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.029) 2.27 ng/u1

response 14749

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	96.29
41.00	42.20	45.22
42.00	29.40	33.41

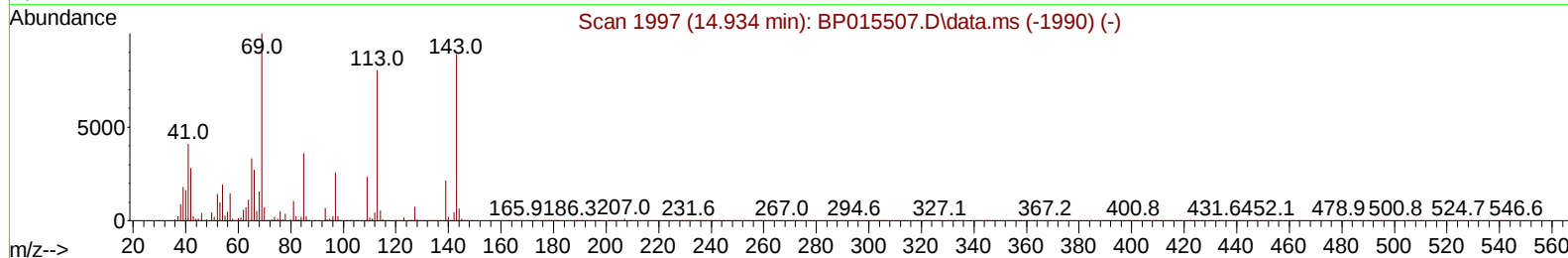
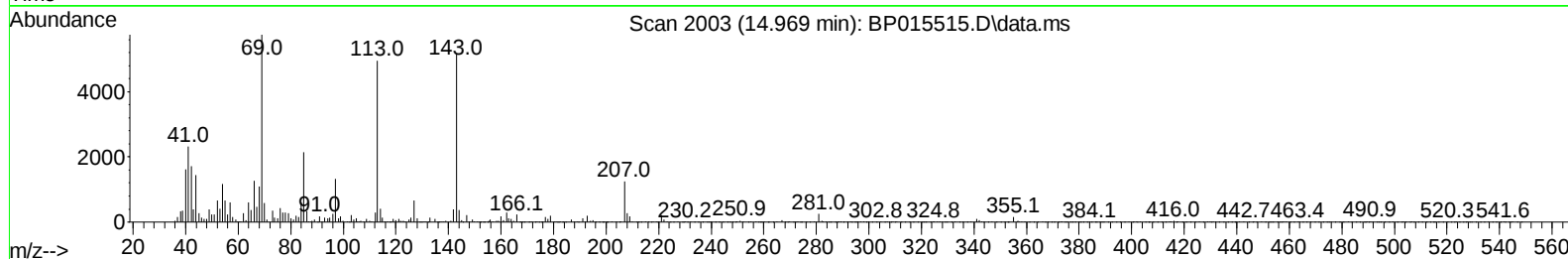
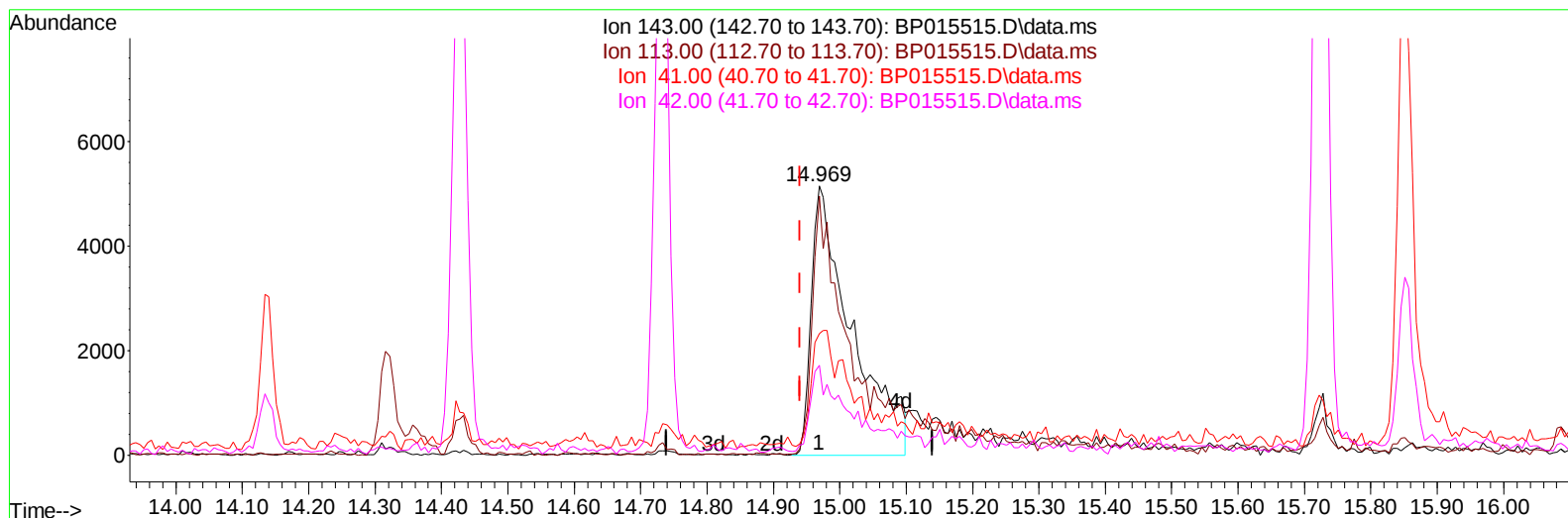
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 Operator : MA/JU
 Sample : 03098-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZEP7

Manual IntegrationsAPPROVED

Quant Time: Jun 14 23:29:55 2023
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 QLast Update : Fri Jun 09 23:33:38 2023
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Reviewed By :Yogesh Patel 06/15/2023
 Supervised By :mohammad ahmed 06/16/2023



TIC: BP015515.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.029) 3.33 ng/ul m

response 21615

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	96.29
41.00	42.20	45.22
42.00	29.40	33.41

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/15/2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	141470	20.000	ng/ul	0.00
20) Naphthalene-d8	10.916	136	652207	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.734	164	459490	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.486	188	1085981	20.000	ng/ul	0.00
79) Chrysene-d12	21.568	240	1203909	20.000	ng/ul	0.00
88) Perylene-d12	24.115	264	1305760	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	17128	4.483	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.246	99	73153	5.613	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.410	67	216160	27.771	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132	202344	21.414	ng/ul	0.00
15) 4-Methylphenol-d8	8.804	113	143588	13.424	ng/ul	0.00
21) Nitrobenzene-d5	9.269	128	139418	27.228	ng/ul	0.00
24) 2-Nitrophenol-d4	9.998	143	152159	26.021	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.540	165	247695	23.221	ng/ul	0.00
31) 4-Chloroaniline-d4	11.098	131	4458m	0.291	ng/ul	0.04
46) Dimethylphthalate-d6	14.139	166	1052929	29.059	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	1109079	27.993	ng/ul	0.00
54) 4-Nitrophenol-d4	14.969	143	21615m	3.331	ng/ul	0.03
60) Fluorene-d10	15.728	176	886009	28.997	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	175711	25.540	ng/ul	0.00
73) Anthracene-d10	17.586	188	1519343	30.749	ng/ul	0.00
81) Pyrene-d10	19.810	212	2097603	32.554	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	2134270	32.576	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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